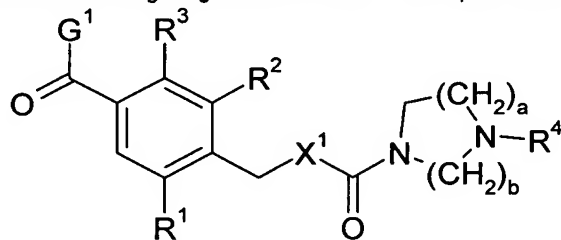


Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

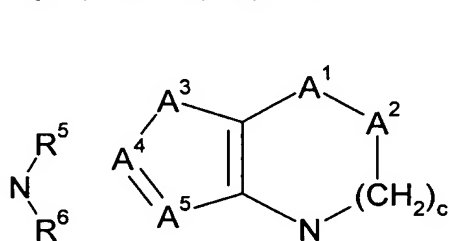
1. (Original) A compound according to general formula 1, or a pharmaceutically acceptable salt thereof



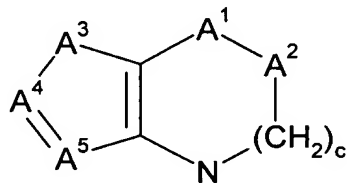
1

wherein:

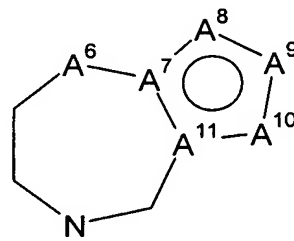
- G¹ is selected from a group according to general formula 2, a group according to general formula 3, a group according to general formula 4, a group according to general formula 5, a group according to general formula 6 and a group according to general formula 7;



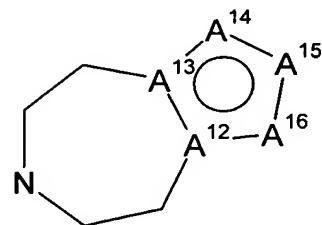
2



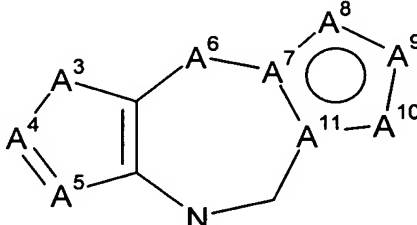
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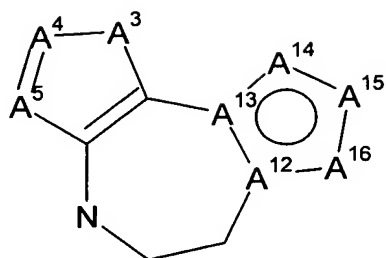
4



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- A¹ is selected from CH₂, CH(OH), NH, N-alkyl, O and S;
- A² is selected from CH₂, CH(OH), C(=O) and NH;
- A³ is selected from S, NH, N-alkyl, -CH=CH- and -CH=N-;
- A⁴ and A⁵ are each selected from CH and N;
- A⁶ is selected from CH₂, NH, N-alkyl and O;
- A⁷ and A¹¹ are selected from C and N;
- A⁸ and A⁹ are selected from CH, N, NH, N(CH₂)_dR⁷ and S;
- A¹⁰ is selected from -CH=CH-, CH, N, NH, N(CH₂)_dR⁷ and S;
- A¹² and A¹³ are selected from N and C;
- A¹⁴, A¹⁵ and A¹⁶ are selected from NH, N-CH₃, S, N and CH;
- X¹ is selected from O and NH;
- R¹, R² and R³ are each selected from H, alkyl, O-alkyl, F, Cl and Br;
- R⁴ is selected from H, alkyl, alkenyl, alkynyl, optionally substituted phenyl, optionally substituted thienyl, optionally substituted furyl, optionally substituted pyridyl, optionally substituted pyrrolyl, optionally substituted pyrazolyl, optionally substituted imidazolyl, optionally substituted oxazolyl, optionally substituted isoxazolyl, optionally substituted thiazolyl, optionally substituted isothiazolyl, -(CH₂)_eR⁸, -CH₂-CH=CH-CH₂-R⁸, -CH₂-C≡C-

CH₂-R⁸, -(CH₂)_g-CH(OH)-(CH₂)_h-R⁸, -(CH₂)_i-O-(CH₂)_j-R⁸ and $\text{CH}_2-\triangle-\text{R}^8$;

- R⁵ and R⁶ are independently selected from alkyl, Ar and -(CH₂)_f-Ar;
- R⁷ is selected from H, alkyl, optionally substituted phenyl, F, OH, O-alkyl, O-acyl, S-alkyl, NH₂, NH-alkyl, N(alkyl)₂, NH-acyl, N(alkyl)-acyl, CO₂H, CO₂-alkyl, CONH₂, CONH-alkyl, CON(alkyl)₂, CN and CF₃;
- R⁸ is selected from H, alkyl, alkenyl, alkynyl, acyl, optionally substituted phenyl, optionally substituted pyridyl, optionally substituted thienyl, optionally substituted furyl, optionally substituted pyrrolyl, optionally substituted pyrazolyl, optionally substituted imidazolyl, optionally substituted oxazolyl, optionally substituted isoxazolyl, optionally substituted thiazolyl, optionally substituted isothiazolyl, F, OH, hydroxyalkyl, O-alkyl, O-acyl, S-alkyl, NH₂, NH-alkyl, N(alkyl)₂, 1-pyrrolidinyl, 1-piperidinyl, 4-morpholinyl, NH-acyl, N(alkyl)-acyl, N₃, CO₂H, CO₂-alkyl, CONH₂, CONH-alkyl, CON(alkyl)₂, CN and CF₃;
- Ar is selected from optionally substituted thienyl and optionally substituted phenyl;

- a is 1 or 2, b is 1, 2 or 3; c is 1 or 2, d is 1, 2 or 3; e is 1, 2, 3 or 4; f is 1, 2 or 3 and g, h, i and j are all independently 1 or 2;

provided that:

- not more than one of A⁸, A⁹ and A¹⁰ is NH, N(CH₂)_dR⁷ or S;
- A⁷ and A¹¹ are not both simultaneously N;
- neither A⁷ nor A¹¹ is N if one of A⁸, A⁹ and A¹⁰ is NH, N(CH₂)_dR⁷ or S;
- if A¹⁰ is not —CH=CH— then one of A⁸, A⁹ and A¹⁰ is NH, N(CH₂)_dR⁷ or S or one of A⁷ and A¹¹ is N;
- not more than one of A¹⁴, A¹⁵ and A¹⁶ is NH, N—CH₃ or S;
- A¹² and A¹³ are not both simultaneously N;
- if one of A¹⁴, A¹⁵ and A¹⁶ is NH, N—CH₃ or S then A¹² and A¹³ are both C; and
- one of A¹⁴, A¹⁵ and A¹⁶ is NH, N—CH₃ or S or one of A¹² and A¹³ is N,

wherein said compound is selected from the group consisting of:

- 4-(3,3-Dimethyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-(2-Cyclopropyl-ethyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-(2-Hydroxymethyl-cyclopropylmethyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-(3-Methyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclopentylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclohexylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;

- 4-(2-Cyclopropyl-ethyl)-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Pentyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Hexyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- (R)-4-(2-Methyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-(2-Ethyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-(2-Methyl-but-2-enyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-ethyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide;
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide; and
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-methoxy-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide.

2. (Original) A pharmaceutical composition comprising a compound according to claim 1 as an active agent.

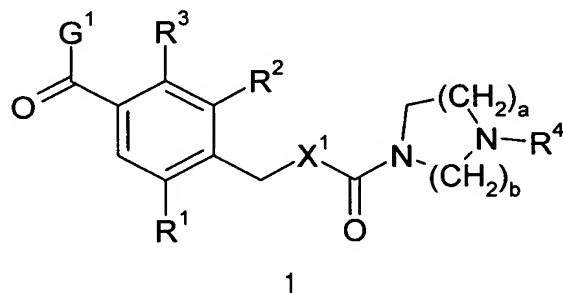
3. (Original) A pharmaceutical composition according to Claim 2 formulated as a tablet or capsule for oral administration.

4. (Currently Amended) A pharmaceutical composition according to Claim 2 ~~or~~ 3 for treatment of primary dysmenorrhoea.

5. (Currently Amended) A pharmaceutical composition according to Claim 2 ~~or~~ 3 for treatment of male erectile dysfunction.

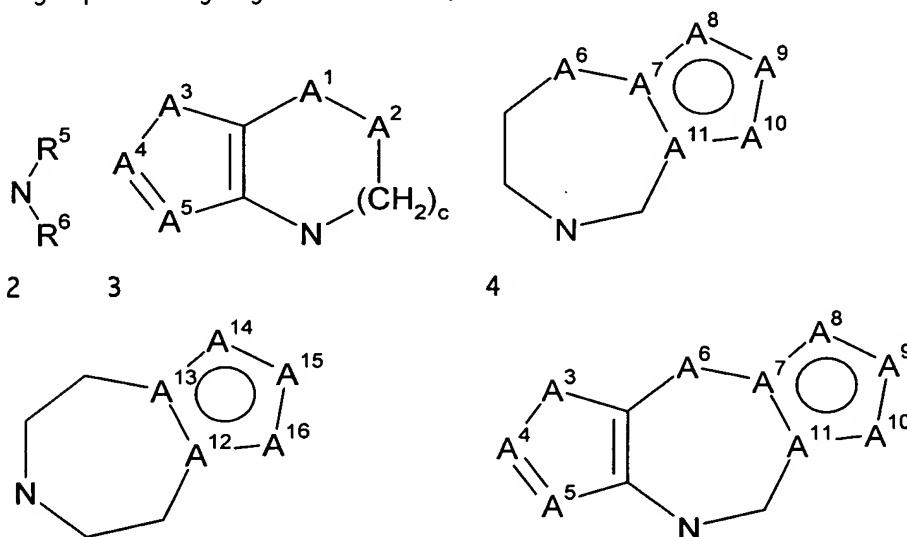
6. (Currently Amended) A pharmaceutical composition according to Claim 2 ~~or~~ 3 for treatment of pre-term labour, hypertension, Raynaud's disease, brain oedema, motion sickness, small cell lung cancer, depression, anxiety, hyponatremia, liver cirrhosis or congestive heart failure.

7. (Original) The use of a compound according to general formula 1, or a pharmaceutically acceptable salt thereof



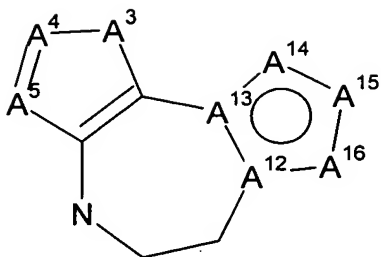
wherein:

- G¹ is selected from a group according to general formula 2, a group according to general formula 3, a group according to general formula 4, a group according to general formula 5, a group according to general formula 6 and a group according to general formula 7;



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- A¹ is selected from CH₂, CH(OH), NH, N-alkyl, O and S;
- A² is selected from CH₂, CH(OH), C(=O) and NH;
- A³ is selected from S, NH, N-alkyl, -CH=CH- and -CH=N-;
- A⁴ and A⁵ are each selected from CH and N;
- A⁶ is selected from CH₂, NH, N-alkyl and O;
- A⁷ and A¹¹ are selected from C and N;
- A⁸ and A⁹ are selected from CH, N, NH, N(CH₂)_dR⁷ and S;
- A¹⁰ is selected from -CH=CH-, CH, N, NH, N(CH₂)_dR⁷ and S;
- A¹² and A¹³ are selected from N and C;
- A¹⁴, A¹⁵ and A¹⁶ are selected from NH, N-CH₃, S, N and CH;
- X¹ is selected from O and NH;
- R¹, R² and R³ are each selected from H, alkyl, O-alkyl, F, Cl and Br;
- R⁴ is selected from H, alkyl, alkenyl, alkynyl, optionally substituted phenyl, optionally substituted thienyl, optionally substituted furyl, optionally substituted pyridyl, optionally substituted pyrrolyl, optionally substituted pyrazolyl, optionally substituted imidazolyl, optionally substituted oxazolyl, optionally substituted isoxazolyl, optionally substituted thiazolyl, optionally substituted isothiazolyl, -(CH₂)_eR⁸, -CH₂-CH=CH-CH₂-R⁸, -CH₂-C≡C-

CH₂-R⁸, -(CH₂)_g-CH(OH)-(CH₂)_h-R⁸, -(CH₂)_i-O-(CH₂)_j-R⁸ and $\text{CH}_2-\triangle-\text{R}^8$;

- R⁵ and R⁶ are independently selected from alkyl, Ar and -(CH₂)_f-Ar;
- R⁷ is selected from H, alkyl, optionally substituted phenyl, F, OH, O-alkyl, O-acyl, S-alkyl, NH₂, NH-alkyl, N(alkyl)₂, NH-acyl, N(alkyl)-acyl, CO₂H, CO₂-alkyl, CONH₂, CONH-alkyl, CON(alkyl)₂, CN and CF₃;
- R⁸ is selected from H, alkyl, alkenyl, alkynyl, acyl, optionally substituted phenyl, optionally substituted pyridyl, optionally substituted thienyl, optionally substituted furyl, optionally substituted pyrrolyl, optionally substituted pyrazolyl, optionally substituted imidazolyl, optionally substituted oxazolyl, optionally substituted isoxazolyl, optionally substituted thiazolyl, optionally substituted isothiazolyl, F, OH, hydroxyalkyl, O-alkyl, O-acyl, S-alkyl,

NH₂, NH-alkyl, N(alkyl)₂, 1-pyrrolidinyl, 1-piperidinyl, 4-morpholinyl, NH-acyl, N(alkyl)-acyl, N₃, CO₂H, CO₂-alkyl, CONH₂, CONH-alkyl, CON(alkyl)₂, CN and CF₃;

- Ar is selected from optionally substituted thienyl and optionally substituted phenyl;

- a is 1 or 2, b is 1, 2 or 3; c is 1 or 2, d is 1, 2 or 3; e is 1, 2, 3 or 4; f is 1, 2 or 3 and g, h, i and j are all independently 1 or 2;

provided that:

- not more than one of A⁸, A⁹ and A¹⁰ is NH, N(CH₂)_dR⁷ or S;

- A⁷ and A¹¹ are not both simultaneously N;

- neither A⁷ nor A¹¹ is N if one of A⁸, A⁹ and A¹⁰ is NH, N(CH₂)_dR⁷ or S;

- if A¹⁰ is not -CH=CH- then one of A⁸, A⁹ and A¹⁰ is NH, N(CH₂)_dR⁷ or S or one of A⁷ and A¹¹ is N;

- not more than one of A¹⁴, A¹⁵ and A¹⁶ is NH, N-CH₃ or S;

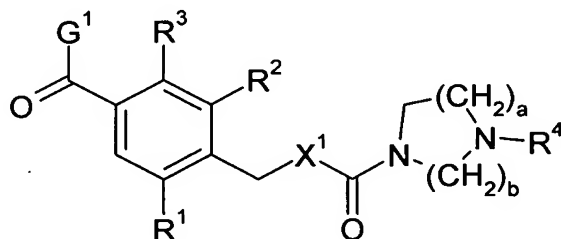
- A¹² and A¹³ are not both simultaneously N;

- if one of A¹⁴, A¹⁵ and A¹⁶ is NH, N-CH₃ or S then A¹² and A¹³ are both C; and

- one of A¹⁴, A¹⁵ and A¹⁶ is NH, N-CH₃ or S or one of A¹² and A¹³ is N,

for the manufacture of a pharmaceutical composition for treatment of primary dysmenorrhoea.

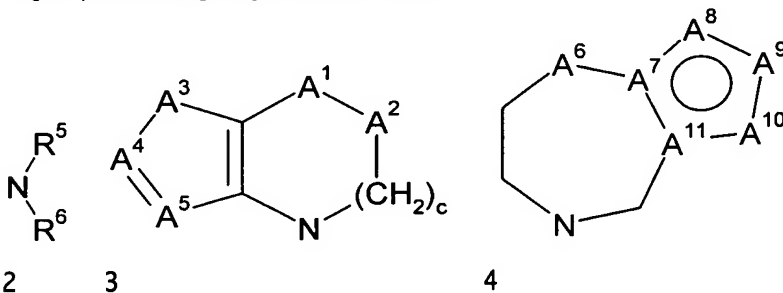
8. (Original) The use of a compound according to general formula 1, or a pharmaceutically acceptable salt thereof

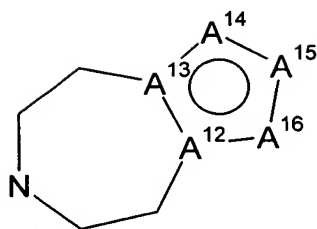


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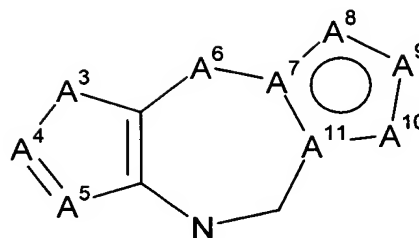
wherein:

- G¹ is selected from a group according to general formula 2, a group according to general formula 3, a group according to general formula 4, a group according to general formula 5, a group according to general formula 6 and a group according to general formula 7;

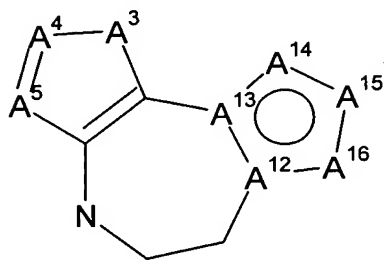




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- A¹ is selected from CH₂, CH(OH), NH, N-alkyl, O and S;
- A² is selected from CH₂, CH(OH), C(=O) and NH;
- A³ is selected from S, NH, N-alkyl, -CH=CH- and -CH=N-;
- A⁴ and A⁵ are each selected from CH and N;
- A⁶ is selected from CH₂, NH, N-alkyl and O;
- A⁷ and A¹¹ are selected from C and N;
- A⁸ and A⁹ are selected from CH, N, NH, N(CH₂)_dR⁷ and S;
- A¹⁰ is selected from -CH=CH-, CH, N, NH, N(CH₂)_dR⁷ and S;
- A¹² and A¹³ are selected from N and C;
- A¹⁴, A¹⁵ and A¹⁶ are selected from NH, N-CH₃, S, N and CH;
- X¹ is selected from O and NH;
- R¹, R² and R³ are each selected from H, alkyl, O-alkyl, F, Cl and Br;
- R⁴ is selected from H, alkyl, alkenyl, alkynyl, optionally substituted phenyl, optionally substituted thienyl, optionally substituted furyl, optionally substituted pyridyl, optionally substituted pyrrolyl, optionally substituted pyrazolyl, optionally substituted imidazolyl, optionally substituted oxazolyl, optionally substituted isoxazolyl, optionally substituted thiazolyl, optionally substituted isothiazolyl, -(CH₂)_eR⁸, -CH₂-CH=CH-CH₂-R⁸, -CH₂-C≡C-CH₂-R⁸, -(CH₂)₉-CH(OH)-(CH₂)_h-R⁸, -(CH₂)_i-O-(CH₂)_j-R⁸ and $\text{CH}_2\text{---}\triangle\text{---R}^8$;
- R⁵ and R⁶ are independently selected from alkyl, Ar and -(CH₂)_f-Ar;

- R⁷ is selected from H, alkyl, optionally substituted phenyl, F, OH, O-alkyl, O-acyl, S-alkyl, NH₂, NH-alkyl, N(alkyl)₂, NH-acyl, N(alkyl)-acyl, CO₂H, CO₂-alkyl, CONH₂, CONH-alkyl, CON(alkyl)₂, CN and CF₃;
- R⁸ is selected from H, alkyl, alkenyl, alkynyl, acyl, optionally substituted phenyl, optionally substituted pyridyl, optionally substituted thienyl, optionally substituted furyl, optionally substituted pyrrolyl, optionally substituted pyrazolyl, optionally substituted imidazolyl, optionally substituted oxazolyl, optionally substituted isoxazolyl, optionally substituted thiazolyl, optionally substituted isothiazolyl, F, OH, hydroxyalkyl, O-alkyl, O-acyl, S-alkyl, NH₂, NH-alkyl, N(alkyl)₂, 1-pyrrolidinyl, 1-piperidinyl, 4-morpholinyl, NH-acyl, N(alkyl)-acyl, N₃, CO₂H, CO₂-alkyl, CONH₂, CONH-alkyl, CON(alkyl)₂, CN and CF₃;
- Ar is selected from optionally substituted thienyl and optionally substituted phenyl;
- a is 1 or 2, b is 1, 2 or 3; c is 1 or 2, d is 1, 2 or 3; e is 1, 2, 3 or 4; f is 1, 2 or 3 and g, h, i and j are all independently 1 or 2;

provided that:

- not more than one of A⁸, A⁹ and A¹⁰ is NH, N(CH₂)_dR⁷ or S;
- A⁷ and A¹¹ are not both simultaneously N;
- neither A⁷ nor A¹¹ is N if one of A⁸, A⁹ and A¹⁰ is NH, N(CH₂)_dR⁷ or S;
- if A¹⁰ is not -CH=CH- then one of A⁸, A⁹ and A¹⁰ is NH, N(CH₂)_dR⁷ or S or one of A⁷ and A¹¹ is N;
- not more than one of A¹⁴, A¹⁵ and A¹⁶ is NH, N-CH₃ or S;
- A¹² and A¹³ are not both simultaneously N;
- if one of A¹⁴, A¹⁵ and A¹⁶ is NH, N-CH₃ or S then A¹² and A¹³ are both C; and
- one of A¹⁴, A¹⁵ and A¹⁶ is NH, N-CH₃ or S or one of A¹² and A¹³ is N,

for the manufacture of a pharmaceutical composition for treatment of pre-term labour, hypertension, Raynaud's disease, brain oedema, motion sickness, small cell lung cancer, depression, anxiety, hyponatremia, liver cirrhosis or congestive heart failure.

9. (Currently Amended) The use according to Claim 7 ~~or 8~~, wherein at least one of R¹, R² and R³ is H and at least one is not H.

10. (Currently Amended) The use according to ~~any one of Claims 7—9~~ Claim 7, wherein one of R¹, R² and R³ is selected from an alkyl group, an O-alkyl group, F, Cl and Br and the others are H.

11. (Currently Amended) The use according to ~~any one of the Claims 7—10~~ Claim 7, wherein X¹ is NH.

12. (Currently Amended) The use according to ~~any one of the Claims 7—11~~ Claim 7, wherein a is 1 and b is 2.

13. (Currently Amended) The use according to ~~any one of the Claims 7—12~~ Claim 7, wherein G¹ is a group according to general formula 3.

14. (Original) The use according to Claim 13, wherein c is 2.

15. (Currently Amended) The use according to Claim 13 ~~or 14~~, wherein A¹ is CH₂ and A² is NH.

16. (Currently Amended) The use according to Claim 13 ~~or 14~~, wherein A¹ is NH or N-alkyl and A² is C(=O).

17. (Currently Amended) The use according to Claim 13 ~~or 14~~, wherein A³ is S and A⁴ and A⁵ are both CH.

18. (Currently Amended) The use according to ~~any of Claims 13—17~~ Claim 13, wherein A³ is —CH=CH- and A⁴ and A⁵ are both CH.

19. (Currently Amended) The use according to ~~any of Claims 13—17~~ Claim 13, wherein A³ is —CH=N- and A⁴ and A⁵ are both CH.

20. (Currently Amended) The use according to ~~any of Claims 13—17~~ Claim 13, wherein A³ is —CH=CH-, A⁴ is CH and A⁵ is N.

21. (Currently Amended) The use according to ~~any of Claims 7—12~~ Claim 7, wherein G¹ is a group according to general formula 6 or 7

22. (Original) The use according to Claim 21, wherein A³ is S and A⁴ and A⁵ are both CH.

23. (Original) The use according to Claim 21, wherein A³ is —CH=CH- and A⁴ and A⁵ are both CH.

24. (Original) The use according to Claim 21, wherein A³ is —CH=N- and A⁴ and A⁵ are both CH.

25. (Original) The use according to Claim 21, wherein A³ is —CH=CH-, A⁴ is CH and A⁵ is N.

26. (Currently Amended) The use according to ~~any one to Claims 7—12~~ Claim 7, wherein G¹ is a group according to general formula 4 or 6.

27. (Original) The use according to Claim 26, wherein A⁶ is NH.
28. (Currently Amended) The use according to Claim 26 ~~or 27~~, wherein A⁸ is NH or N-(CH₂)_d-R⁷.
29. (Original) The use according to Claim 28, wherein A⁹ is N and A¹⁰ is CH.
30. (Currently Amended) The use according to Claim 7 ~~or 8~~, wherein one of R¹, R² and R³ is selected from an alkyl group, an O-alkyl group, F, Cl and Br and the others are H and X¹ is NH.
31. (Currently Amended) The use according to ~~any one of Claims 7 or 8~~ Claim 7, wherein one of R¹, R² and R³ is selected from an alkyl group, an O-alkyl group, F, Cl and Br and the others are H and X¹ is NH, a is 1 and b is 2.
32. (Currently Amended) The use according to Claim 7 ~~or 8~~, wherein G¹ is a group according to general formula 6, A⁴, A⁵ and A¹⁰ are all CH, A⁶ is NH, A⁷ and A¹¹ are both C, A⁸ is N-(CH₂)_d-R⁷ and A⁹ is N.
33. (Currently Amended) The use according to Claim 7 ~~or 8~~, wherein R¹ is an alkyl group, an O-alkyl group, F, Cl or Br, R² and R³ are both H, X¹ is NH, a is 1, b is 2, G¹ is a group according to general formula 6, A⁴, A⁵ and A¹⁰ are all CH, A⁶ is NH, A⁷ and A¹¹ are both C, A⁸ is N-(CH₂)_d-R⁷ and A⁹ is N.
34. (Currently Amended) The use according to Claim 7 ~~or 8~~, wherein said compound is selected from the group consisting of:
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-(3,3-Dimethyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-(3-Methylsulfanyl-propyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-(2-Cyclopropyl-ethyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,

- 4-(2-Hydroxymethyl-cyclopropylmethyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(3-Methyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopentylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclohexylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(2-Cyclopropyl-ethyl)-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Pentyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Hexyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- (R)-4-(2-Methyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(2-Ethyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(2-Methyl-but-2-enyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,

- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-ethyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide, and
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-methoxy-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide.

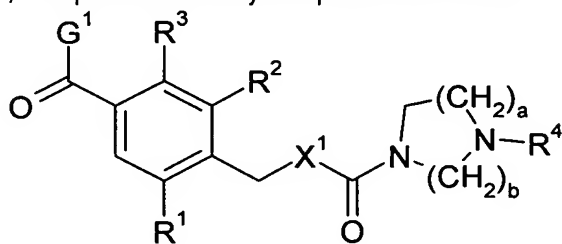
35. (Original) The use of a compound selected from the group consisting of:

- 4-(3,3-Dimethyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(2-Cyclopropyl-ethyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(2-Hydroxymethyl-cyclopropylmethyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(3-Methyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopentylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclohexylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,

- 4-(2-Cyclopropyl-ethyl)-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Pentyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Hexyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - (R)-4-(2-Methyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-(2-Ethyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-(2-Methyl-but-2-enyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-ethyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide, and
 - 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-methoxy-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- for the manufacture of a pharmaceutical composition for treatment of male erectile dysfunction.

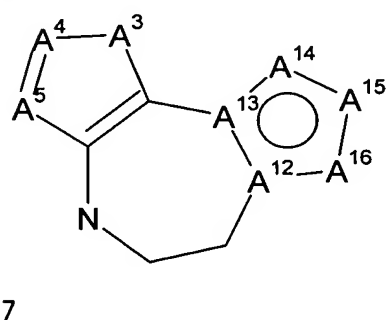
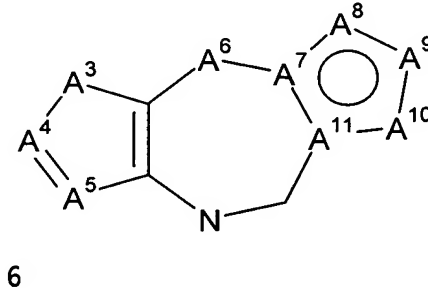
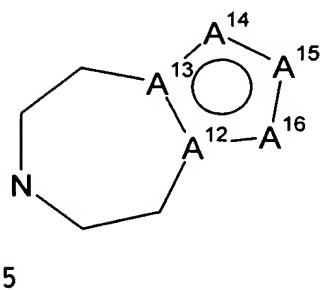
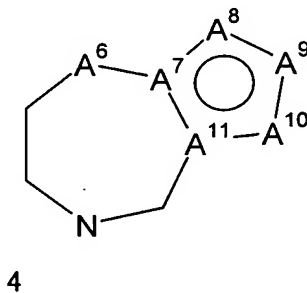
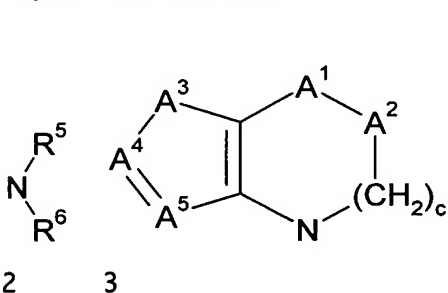
36. (Original) A method for treatment of a disorder selected from the group consisting of primary dysmenorrhoea, pre-term labour, hypertension, Raynaud's disease, brain oedema, motion sickness, small cell lung cancer, depression, anxiety, hyponatremia, liver cirrhosis and congestive heart failure which comprises the

administration to a person in need of such treatment of therapeutically effective amount of a compound according to general formula 1, or a pharmaceutically acceptable salt thereof



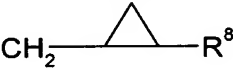
wherein:

- G¹ is selected from a group according to general formula 2, a group according to general formula 3, a group according to general formula 4, a group according to general formula 5, a group according to general formula 6 and a group according to general formula 7;



- A¹ is selected from CH₂, CH(OH), NH, N-alkyl, O and S;
 - A² is selected from CH₂, CH(OH), C(=O) and NH;
 - A³ is selected from S, NH, N-alkyl, -CH=CH- and -CH=N-;

- A⁴ and A⁵ are each selected from CH and N;
- A⁶ is selected from CH₂, NH, N-alkyl and O;
- A⁷ and A¹¹ are selected from C and N;
- A⁸ and A⁹ are selected from CH, N, NH, N(CH₂)_dR⁷ and S;
- A¹⁰ is selected from -CH=CH-, CH, N, NH, N(CH₂)_dR⁷ and S;
- A¹² and A¹³ are selected from N and C;
- A¹⁴, A¹⁵ and A¹⁶ are selected from NH, N-CH₃, S, N and CH;
- X¹ is selected from O and NH;
- R¹, R² and R³ are each selected from H, alkyl, O-alkyl, F, Cl and Br;
- R⁴ is selected from H, alkyl, alkenyl, alkynyl, optionally substituted phenyl, optionally substituted thienyl, optionally substituted furyl, optionally substituted pyridyl, optionally substituted pyrrolyl, optionally substituted pyrazolyl, optionally substituted imidazolyl, optionally substituted oxazolyl, optionally substituted isoxazolyl, optionally substituted thiazolyl, optionally substituted isothiazolyl, -(CH₂)_eR⁸, -CH₂-CH=CH-CH₂-R⁸, -CH₂-C≡C-

CH₂-R⁸, -(CH₂)_g-CH(OH)-(CH₂)_h-R⁸, -(CH₂)_i-O-(CH₂)_j-R⁸ and  ;

- R⁵ and R⁶ are independently selected from alkyl, Ar and -(CH₂)_f-Ar;
- R⁷ is selected from H, alkyl, optionally substituted phenyl, F, OH, O-alkyl, O-acyl, S-alkyl, NH₂, NH-alkyl, N(alkyl)₂, NH-acyl, N(alkyl)-acyl, CO₂H, CO₂-alkyl, CONH₂, CONH-alkyl, CON(alkyl)₂, CN and CF₃;
- R⁸ is selected from H, alkyl, alkenyl, alkynyl, acyl, optionally substituted phenyl, optionally substituted pyridyl, optionally substituted thienyl, optionally substituted furyl, optionally substituted pyrrolyl, optionally substituted pyrazolyl, optionally substituted imidazolyl, optionally substituted oxazolyl, optionally substituted isoxazolyl, optionally substituted thiazolyl, optionally substituted isothiazolyl, F, OH, hydroxyalkyl, O-alkyl, O-acyl, S-alkyl, NH₂, NH-alkyl, N(alkyl)₂, 1-pyrrolidinyl, 1-piperidinyl, 4-morpholinyl, NH-acyl, N(alkyl)-acyl, N₃, CO₂H, CO₂-alkyl, CONH₂, CONH-alkyl, CON(alkyl)₂, CN and CF₃;
- Ar is selected from optionally substituted thienyl and optionally substituted phenyl;
- a is 1 or 2, b is 1, 2 or 3; c is 1 or 2, d is 1, 2 or 3; e is 1, 2, 3 or 4; f is 1, 2 or 3 and g, h, i and j are all independently 1 or 2;

provided that:

- not more than one of A⁸, A⁹ and A¹⁰ is NH, N(CH₂)_dR⁷ or S;
- A⁷ and A¹¹ are not both simultaneously N;
- neither A⁷ nor A¹¹ is N if one of A⁸, A⁹ and A¹⁰ is NH, N(CH₂)_dR⁷ or S;
- if A¹⁰ is not -CH=CH- then one of A⁸, A⁹ and A¹⁰ is NH, N(CH₂)_dR⁷ or S or one of A⁷ and A¹¹ is N;
- not more than one of A¹⁴, A¹⁵ and A¹⁶ is NH, N-CH₃ or S;
- A¹² and A¹³ are not both simultaneously N;
- if one of A¹⁴, A¹⁵ and A¹⁶ is NH, N-CH₃ or S then A¹² and A¹³ are both C; and

- one of A¹⁴, A¹⁵ and A¹⁶ is NH, N—CH₃ or S or one of A¹² and A¹³ is N.

37. (Original) The method of Claim 36, wherein at least one of R¹, R² and R³ is H and at least one is not H.

38. (Original) The method of Claim 36, wherein one of R¹, R² and R³ is selected from an alkyl group, an O-alkyl group, F, Cl and Br and the others are H.

39. (Original) The method of Claim 36, wherein X¹ is NH.

40. (Original) The method of Claim 36, wherein a is 1 and b is 2.

41. (Original) The method of Claim 36, wherein G¹ is a group according to general formula 3.

42. (Original) The method of to Claim 41, wherein c is 2.

43. (Original) The method of Claim 41, wherein A¹ is CH₂ and A² is NH.

44. (Original) The method of Claim 41, wherein A¹ is NH or N-alkyl and A² is C(=O).

45. (Original) The method of Claim 41, wherein A³ is S and A⁴ and A⁵ are both CH.

46. (Original) The method of Claim 41, wherein A³ is —CH=CH- and A⁴ and A⁵ are both CH.

47. (Original) The method of Claim 41, wherein A³ is —CH=N- and A⁴ and A⁵ are both CH.

48. (Original) The method of Claim 41, wherein A³ is —CH=CH-, A⁴ is CH and A⁵ is N.

49. (Original) The method of Claim 36, wherein G¹ is a group according to general formula 6 or 7.

50. (Original) The method of Claim 49, wherein A³ is S and A⁴ and A⁵ are both CH.

51. (Original) The method of Claim 49, wherein A³ is —CH=CH- and A⁴ and A⁵ are both CH.

52. (Original) The method of Claim 49, wherein A³ is —CH=N- and A⁴ and A⁵ are both CH.

53. (Original) The method of Claim 49, wherein A³ is —CH=CH—, A⁴ is CH and A⁵ is N.
54. (Original) The method of Claim 36, wherein G¹ is a group according to general formula 4 or 6.
55. (Original) The method of Claim 54, wherein A⁶ is NH.
56. (Original) The method of Claim 54, wherein A⁸ is NH or N-(CH₂)_d-R⁷.
57. (Original) The method of Claim 56, wherein A⁹ is N and A¹⁰ is CH.
58. (Original) The method of Claim 36, wherein one of R¹, R² and R³ is selected from an alkyl group, an O-alkyl group, F, Cl and Br and the others are H and X¹ is NH.
59. (Original) The method of Claim 36, wherein one of R¹, R² and R³ is selected from an alkyl group, an O-alkyl group, F, Cl and Br and the others are H and X¹ is NH, a is 1 and b is 2.
60. (Original) The method of Claim 36, wherein G¹ is a group according to general formula 6, A⁴, A⁵ and A¹⁰ are all CH, A⁶ is NH, A⁷ and A¹¹ are both C, A⁸ is N-(CH₂)_d-R⁷ and A⁹ is N.
61. (Original) The method of Claim 36, wherein R¹ is an alkyl group, an O-alkyl group, F, Cl or Br, R² and R³ are both H, X¹ is NH, a is 1, b is 2, G¹ is a group according to general formula 6, A⁴, A⁵ and A¹⁰ are all CH, A⁶ is NH, A⁷ and A¹¹ are both C, A⁸ is N-(CH₂)_d-R⁷ and A⁹ is N.
62. (Original) The method of Claim 36, wherein said compound is selected from the group consisting of:
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-(3,3-Dimethyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-(3-Methylsulfanyl-propyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-(2-Cyclopropyl-ethyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,

- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(2-Hydroxymethyl-cyclopropylmethyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(3-Methyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopentylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclohexylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(2-Cyclopropyl-ethyl)-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Pentyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Hexyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- (R)-4-(2-Methyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(2-Ethyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(2-Methyl-but-2-enyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,

- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-ethyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide, and
- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-methoxy-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide.

63. (Original) A method for treatment of male erectile dysfunction which comprises the administration to a person in need of such treatment of therapeutically effective amount of a compound selected from the group consisting of:

- 4-(3,3-Dimethyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(2-Cyclopropyl-ethyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(2-Hydroxymethyl-cyclopropylmethyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-(3-Methyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopentylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclohexylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
- 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 3-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,

- 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-(2-Cyclopropyl-ethyl)-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Pentyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Hexyl-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - (R)-4-(2-Methyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-(2-Ethyl-butyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-(2-Methyl-but-2-enyl)-piperazine-1-carboxylic acid 2-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-methyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-fluoro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-ethyl-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 2-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclopropylmethyl-piperazine-1-carboxylic acid 2-chloro-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide,
 - 4-Cyclobutylmethyl-piperazine-1-carboxylic acid 3-methoxy-4-(3-methyl-4,10-dihydro-3H-2,3,4,9-tetraaza-benzo[f]azulene-9-carbonyl)-benzylamide; and
- pharmaceutically acceptable salts of the above mentioned compounds.